

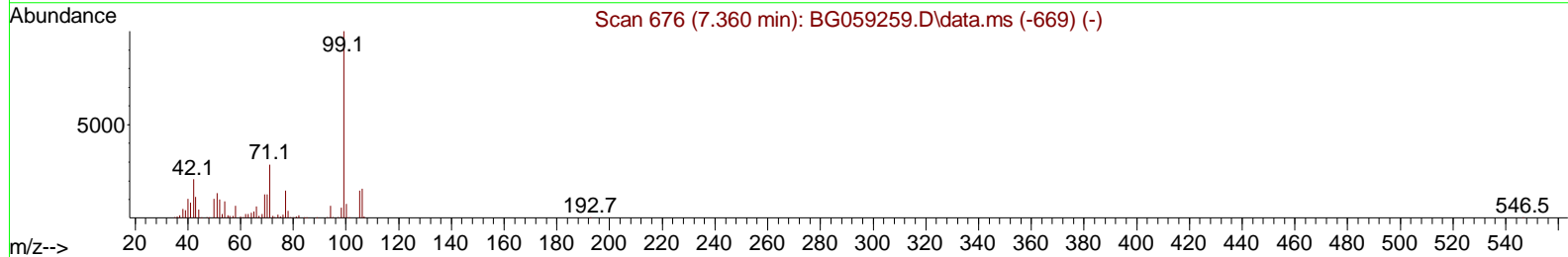
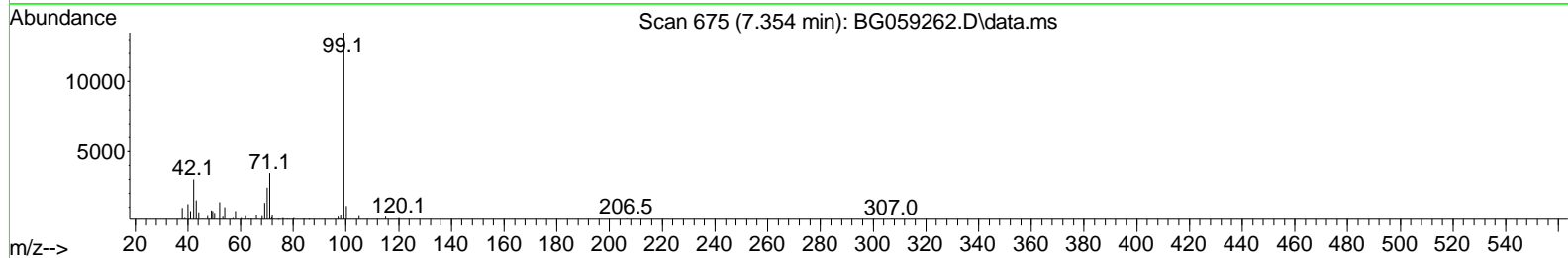
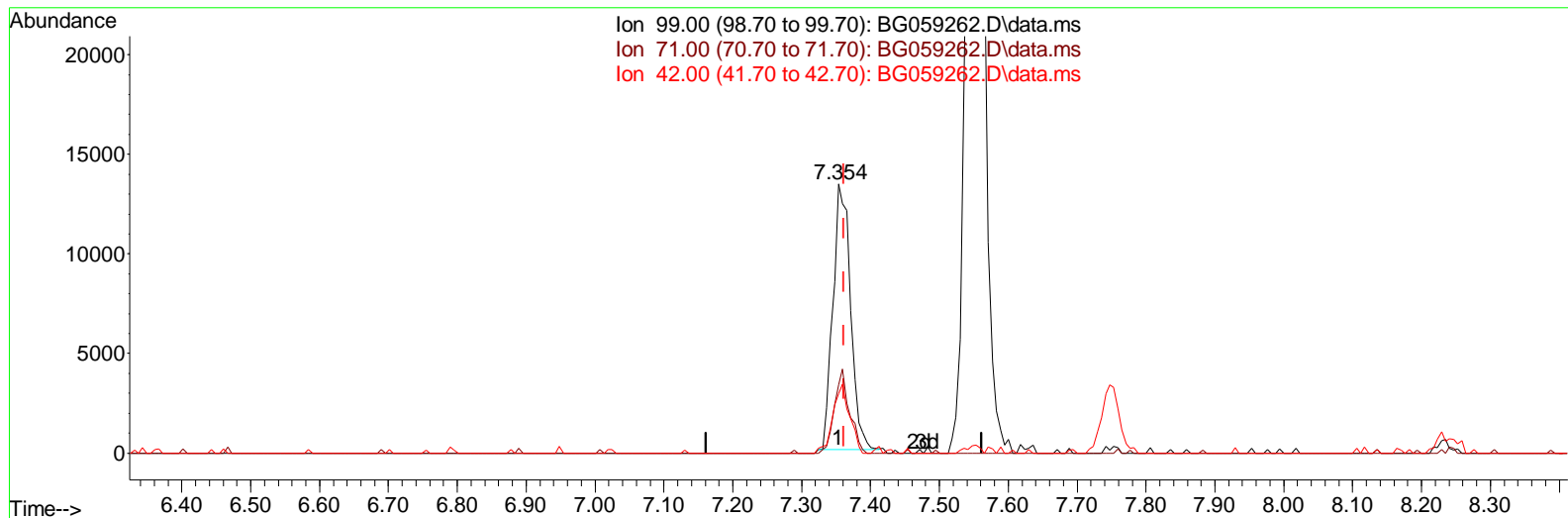
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100323\
 Data File : BG059262.D
 Acq On : 3 Oct 2023 13:09
 Operator : MA/JU
 Sample : 04480-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
ClientSampleId :
 BBJ27

Manual IntegrationsAPPROVED

Quant Time: Oct 03 23:33:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 04:33:25 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/04/2023
 Supervised By :mohammad ahmed 10/04/2023



TIC: BG059262.D\data.ms

(7) Phenol-d5 (S)

7.354min (-0.008) 4.49 ng/ul

response 23648

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	27.00	25.46
42.00	24.00	22.08
0.00	0.00	0.00

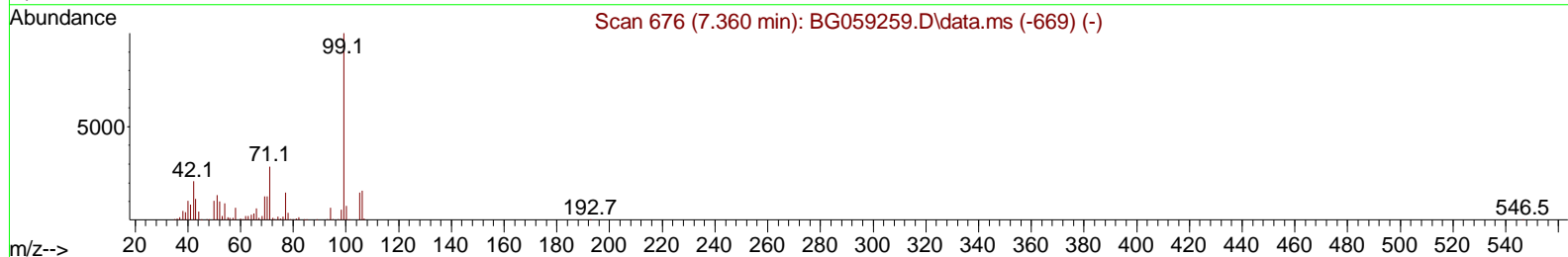
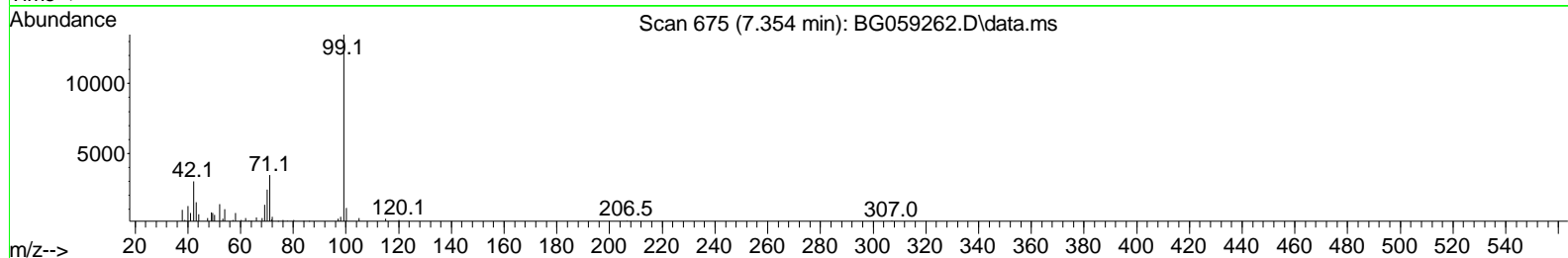
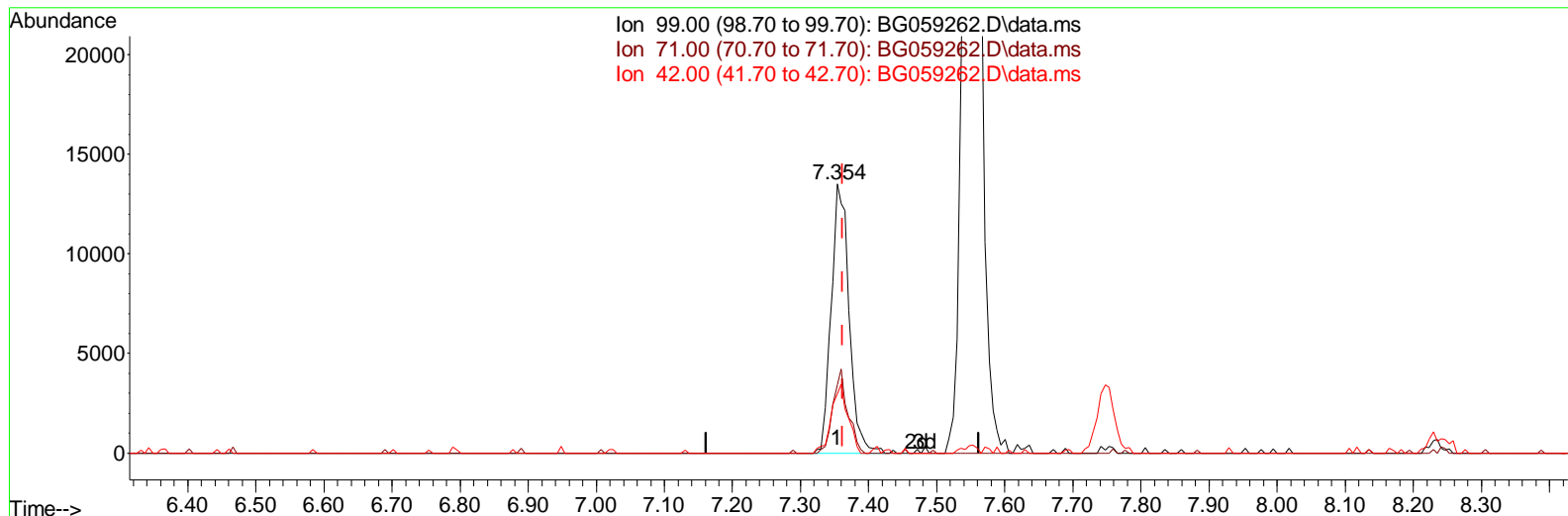
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TIC: BG059262.D\data.ms

(7) Phenol-d5 (S)

7.354min (-0.008) 4.70 ng/ul m

response	24761
Ion	Exp% Act%
99.00	100.00 100.00
71.00	27.00 25.46
42.00	24.00 22.08
0.00	0.00 0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100323\
 Data File : BG059262.D
 Acq On : 3 Oct 2023 13:09
 Operator : MA/JU
 Sample : 04480-02
 Mi sc :
 ALS Vial : 5 Sample Multi plier: 1

Instrument :
 BNA_G
ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/04/2023
 Supervised By :mohammad ahmed 10/04/2023

Quant Time: Oct 03 23: 43: 49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
 Quant Title : SVOA CALI BRATI ON
 QLast Update : Thu Sep 28 04: 33: 25 2023
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
Internal Standards						
1) 1, 4-Di chl orobenzene-d4	8.235	152	52650	20.000	ng/ul	0.00
20) Naphthal ene-d8	11.073	136	259839	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.862	164	165525	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.612	188	371139	20.000	ng/ul	0.00
79) Chrysene-d12	21.913	240	330563	20.000	ng/ul	0.00
88) Perylene-d12	25.391	264	389527	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1, 4-Di oxane-d8	3.535	96	5632	4.037	ng/uL	0.00
4) Pyri di ne-d5	3.969	84	32945	8.094	ng/ul	0.00
7) Phenol -d5	7.354	99	24761m	4.698	ng/ul	0.00
9) Bi s-(2-Chl oroethyl)eth. . .	7.548	67	78550	23.332	ng/ul	0.00
11) 2-Chl orophenol -d4	7.747	132	68084	19.218	ng/ul	0.00
15) 4-Methyl phenol -d8	8.922	113	42872	10.524	ng/ul	0.00
21) Ni trobenzene-d5	9.428	128	46160	25.175	ng/ul	0.00
24) 2-Ni trophenol -d4	10.144	143	47087	26.159	ng/ul	0.00
28) 2, 4-Di chl orophenol -d3	10.673	165	83012	22.450	ng/ul	0.00
31) 4-Chl oroani li ne-d4	11.220	131	80084	13.386	ng/ul	0.00
46) Di methyl phthal ate-d6	14.257	166	297409	24.851	ng/ul	0.00
49) Acenaphthyl ene-d8	14.569	160	351802	23.420	ng/ul	0.00
54) 4-Ni trophenol -d4	15.050	143	7248	3.468	ng/ul	0.00
60) Fl uorene-d10	15.850	176	268847	23.603	ng/ul	0.00
65) 4, 6-Di ni tro-2-methyl ph. . .	15.961	200	38091	19.967	ng/ul	0.00
73) Anthracene-d10	17.712	188	408046	24.129	ng/ul	0.00
81) Pyrene-d10	19.986	212	498472	26.927	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.145	264	475874	25.422	ng/ul	-0.02

Target Compounds Qval ue

(#) = qual i fi er out of range (m) = manual i ntegrati on (+) = signal s summed

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